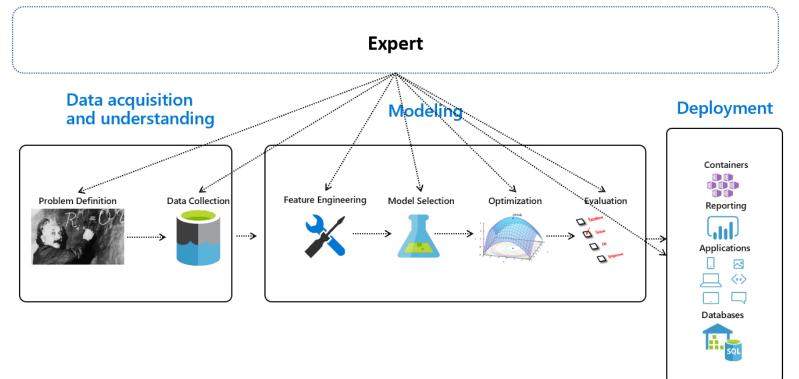
# **Automatic Machine Learning**

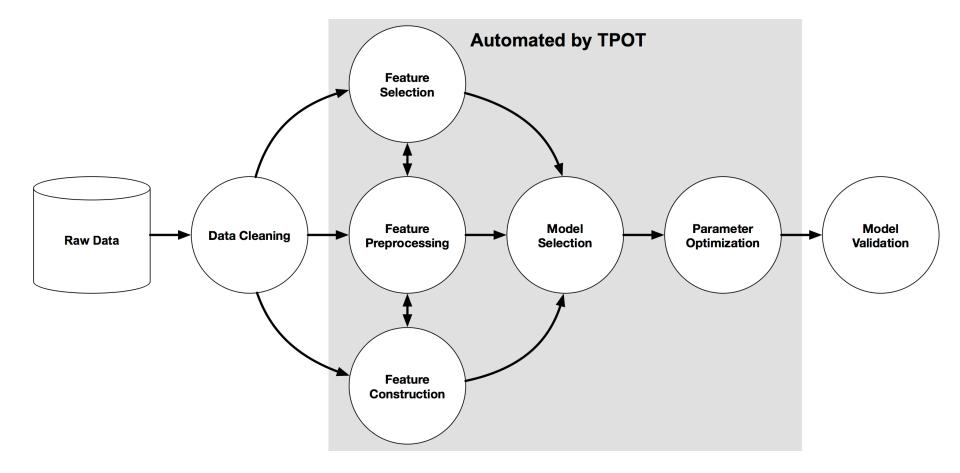
#### Jianping Fan Department of Computer Science UNC-Charlotte

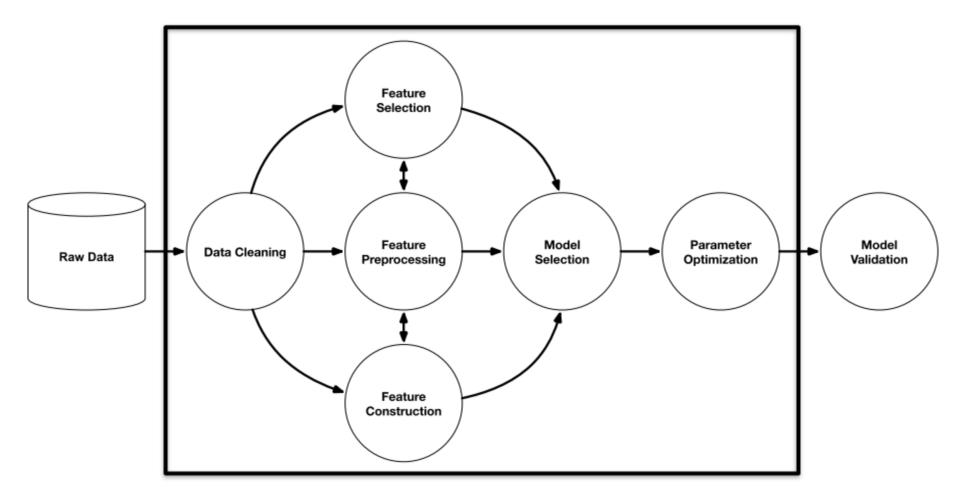
Course Website: http://webpages.uncc.edu/jfan/itcs5152.html

#### **End-to-end Machine Learning Pipeline**

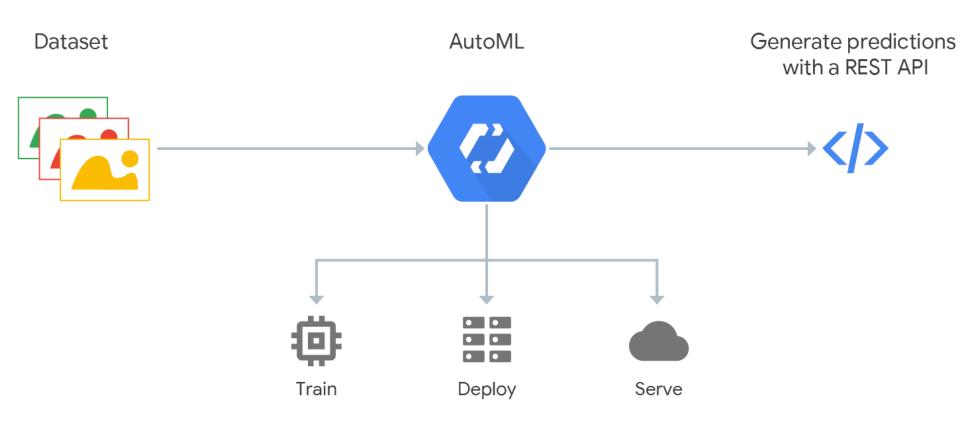








#### Google Cloud AutoML



In the definition of Automated Machine Learning, AutoML includes:

(a) automated feature engineering

(b) automated model selection and hyperparameter tuning

(c) automated neural network architecture selection

(d) automated deployment

The future of AutoML in the next 3–5 years:

- (a) AutoML will also handle most of the data cleaning process
- (b) AutoML will vastly improve deep learning
- (c) AutoML will scaled to large data sets
- (d) AutoML will become human competitive
- (e) AutoML will transform the practice of data science as we know it
- (f) AutoML is only a small part of a greater metalearning movement

## The advantages of AutoML Default parameters are almost always bad



#### AutoML handles this for you!

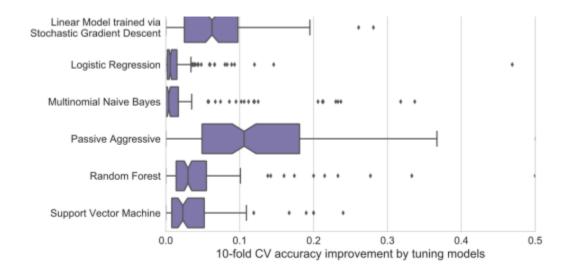


Image source: R. Olson & W. La Cava et. al. (2017) "Data-driven advice for applying machine learning to bioinformatics problems."

## The advantages of AutoML There is no "best" ML algorithm for all problems



% out of 165 datasets where model A outperformed model B

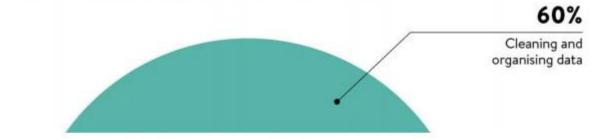
#### AutoML handles this for you!

AdaBoost -	1%	7%	10%	15%	30%		32%			47%	59%	76%	77%
Logistic Regression -	5%	10%	3%	8%	11%	31%	33%	35%			54%	79%	81%
Passive Aggressive -	2%	6%	1%	5%	0%	18%	28%	28%	13%		50%	81%	79%
Bernoulli Naive Bayes -	0%	2%	2%	4%	10%	13%	18%	15%	22%	25%		62%	68%
Gaussian Naive Bayes -	0%	1%	3%	2%	6%	6%	11%	12%	9%	10%	22%		45%
Multinomial Naive Bayes -	1%	1%	2%	2%	2%	5%	10%	14%	4%	5%	13%	39%	
	бтв	ŔF	svм	ERF	sģD	ĸŃN	τα	ÅВ	LR	PA	BNB	GNB	MNB
		Losses											

Image source: R. Olson & W. La Cava et. al. (2017) "Data-driven advice for applying machine learning to bioinformatics problems."

### The advantages of AutoML AutoML is a huge time-saver

WHAT DATA SCIENTISTS SPEND THE MOST TIME DOING



### AutoML handles some of this for you!

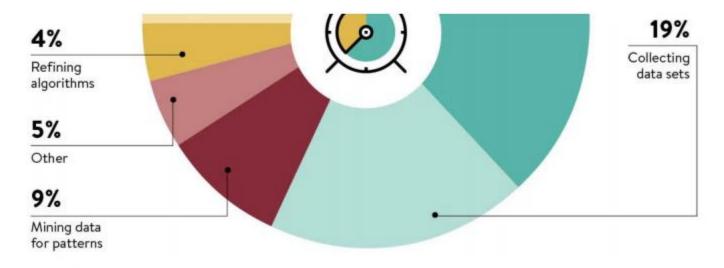


Image source: visit.crowdflower.com/data-science-report.html

### Early AutoML focused on only parameter tuning

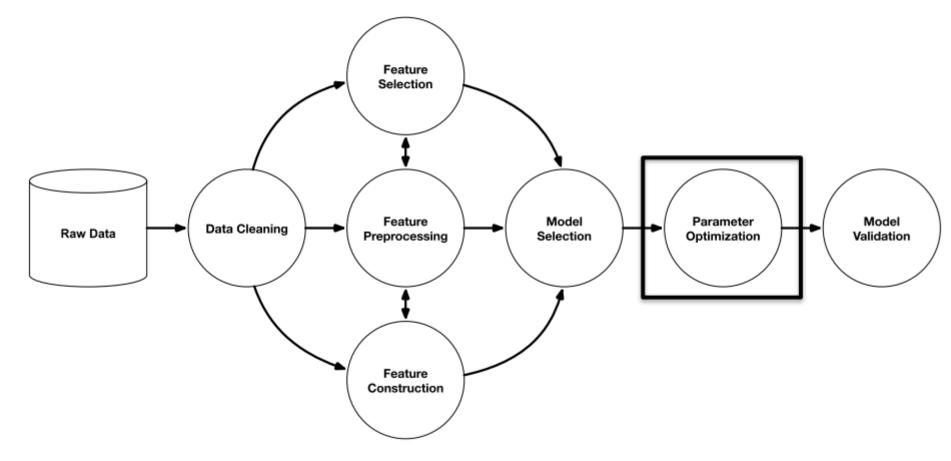
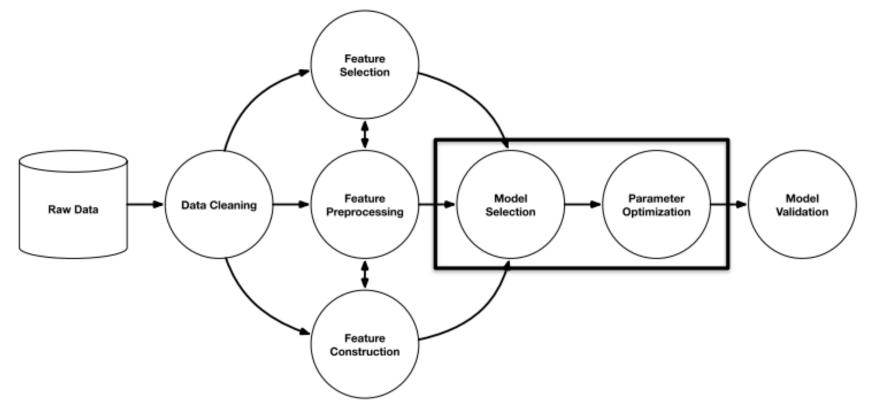


Image source: R. Olson et. al. (2016) "Evaluation of a Tree-based Pipeline Optimization Tool for Automating Data Science."

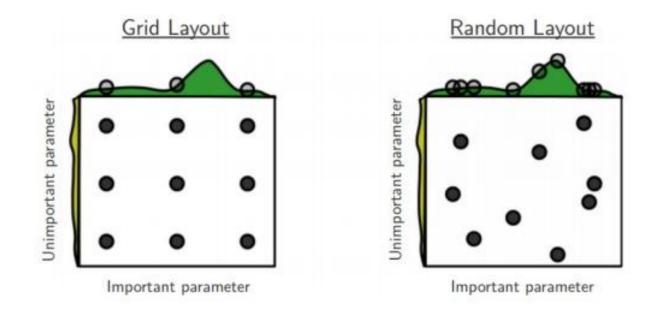
#### Early AutoML focused on only parameter tuning



#### ... and maybe (limited) model selection

Image source: R. Olson et. al. (2016) "Evaluation of a Tree-based Pipeline Optimization Tool for Automating Data Science."

### We mostly used grid search and random search



#### Nowadays, we wouldn't really call this AutoML

#### Modern AutoML optimizes the entire ML workflow

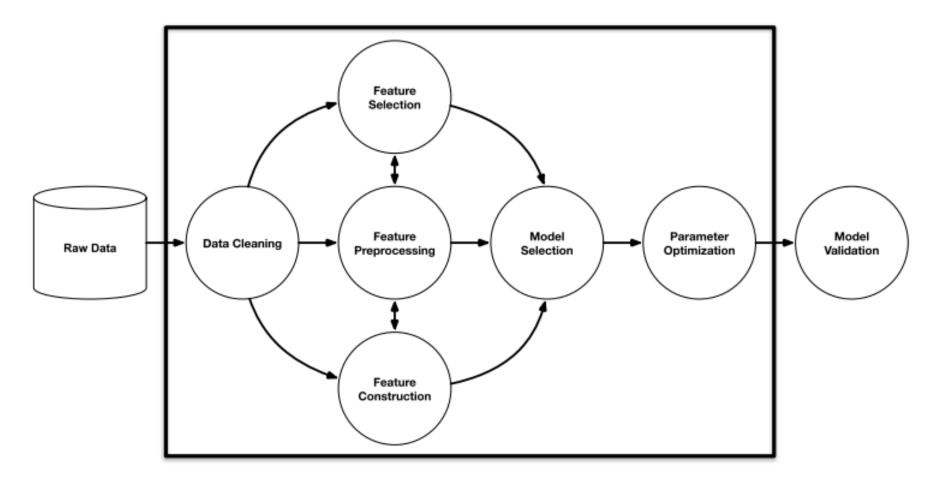
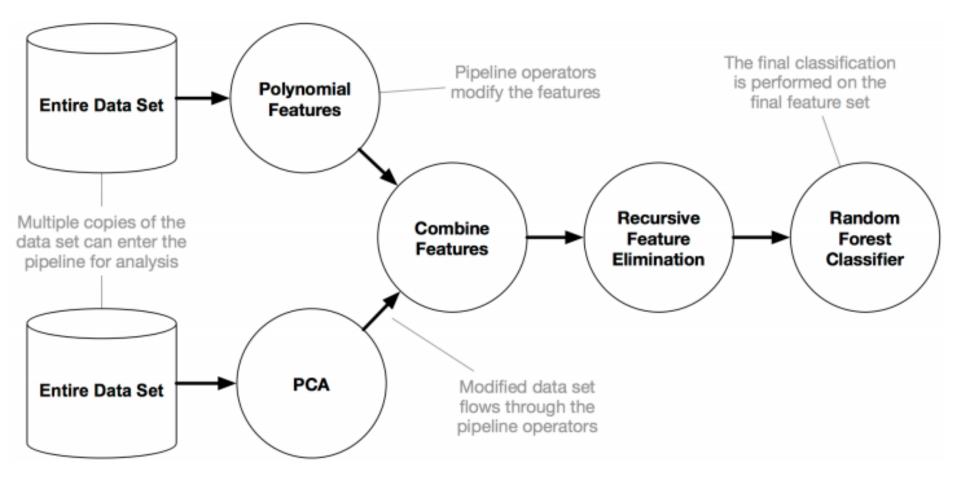


Image source: R. Olson et. al. (2016) "Evaluation of a Tree-based Pipeline Optimization Tool for Automating Data Science."

# Machine Learning analyses can be complex



# Larger search spaces require smarter optimization

- Meta-learning
- Bayesian optimization
- Genetic programming
- Multi-armed bandit

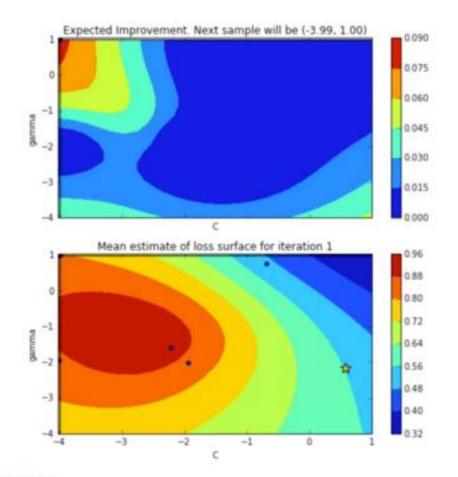


Image source: thuijskens.github.io/2016/12/29/bayesian-optimisation/

### AutoML in the near future

- AutoML will also handle most of the data cleaning process
  - Unstructured data → tabular data ready for analysis
  - Capture & automate human approaches to data cleaning
- AutoML will vastly improve Deep Learning
  - Automated DNN architecture design
  - Automated preprocessing of data prior to modeling

#### AutoML will scale to large datasets

- Most AutoML tools are very slow on "Big Data"
- Spark, Dask, TensorFlow, etc. will help bring AutoML to scale
- H2O AutoML and TPOT on Dask already shows promise for scalability

#### AutoML will become human-competitive

- Already human-competitive on several Kaggle challenges (top 10% of entries)
- Already human-competitive in DNN architecture design (Google AutoML)

#### AutoML in the future

- AutoML will transform the practice of data science as we know it
  - "Data Science Assistant" → Junior Data Scientist level
  - Less focus on choosing the right ML workflow
  - More focus on posing the right questions, collecting & curating the right data, and "thinking like a data scientist"
- AutoML is only a small part of a greater meta-learning movement
  - Computer programming is focused on automating rote tasks
  - Machine learning is focused on automating the automation of rote tasks
  - Meta-learning is focused on automating the automation of automation
    - i.e., enabling the machine to learn how to learn in the best way possible

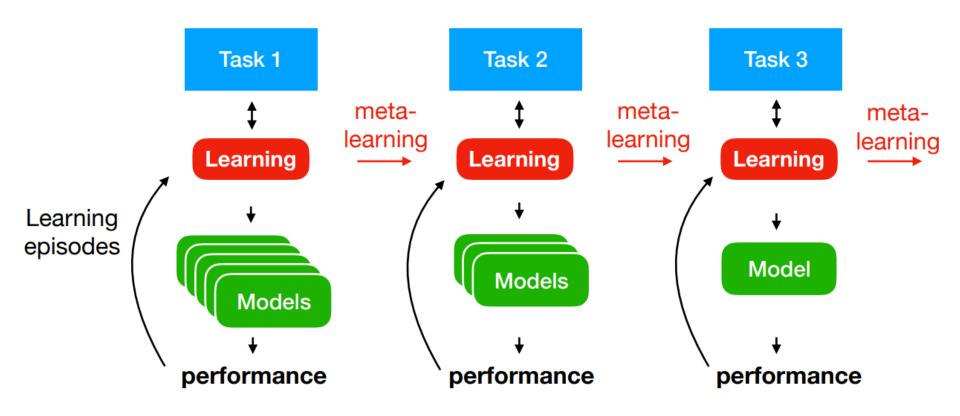
Frank Hutter, Joaquin Vanschoren, NeurIPS 2018 Tutorial on Automatic Machine Learning

#### Learning to Learn

# Learning is a never-ending process

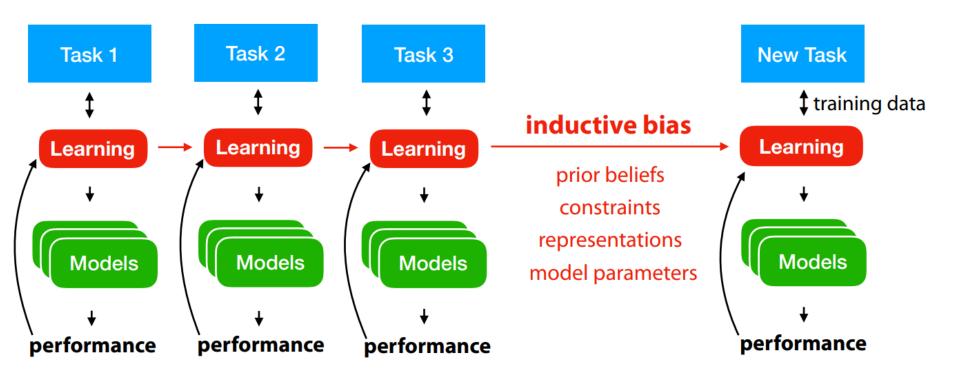
Tasks come and go, but learning is forever

Learn more effectively: less trial-and-error, less data



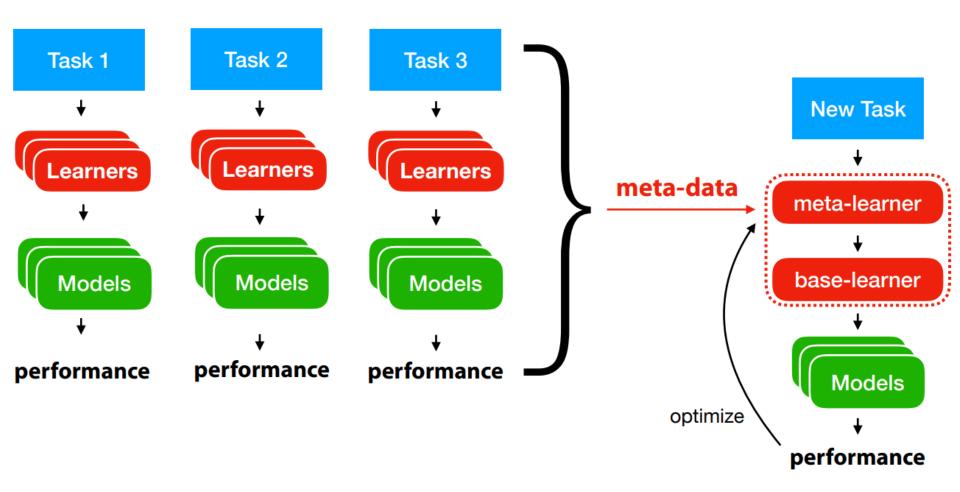
# Learning to learn

**Inductive bias**: all assumptions added to the training data to learn effectively If prior tasks are *similar*, we can *transfer* prior knowledge to new tasks (if not it may actually harm learning)



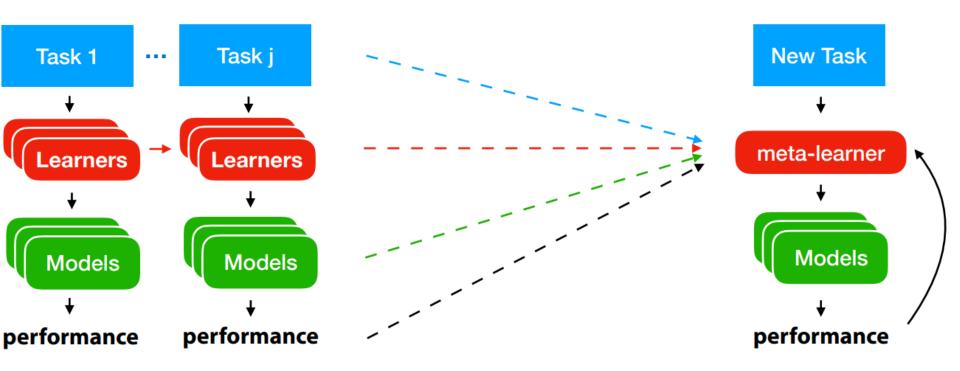
# Meta-learning

Collect meta-data about learning episodes and learn from them Meta-learner *learns* a (base-)learning algorithm, *end-to-end* 



# Three approaches for increasingly similar tasks

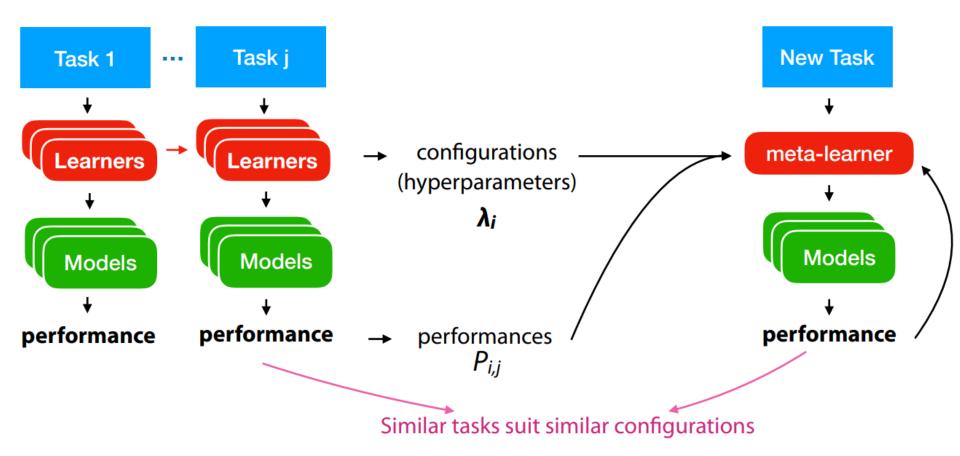
- 1. Transfer prior knowledge about what generally works well
- 2. Reason about model performance across tasks
- 3. Start from models trained earlier on similar tasks



# 1. Learning from prior evaluations

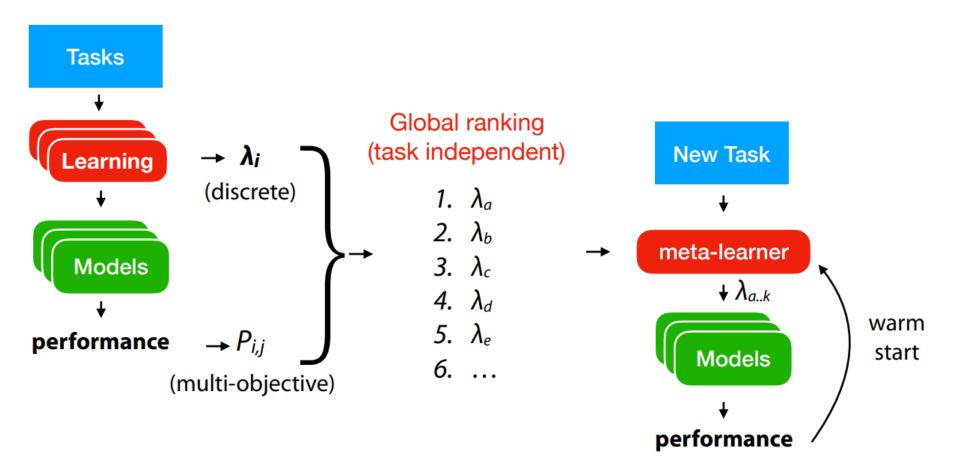
#### Configurations: settings that uniquely define the model

(algorithm, pipeline, neural architecture, hyper-parameters, ...)



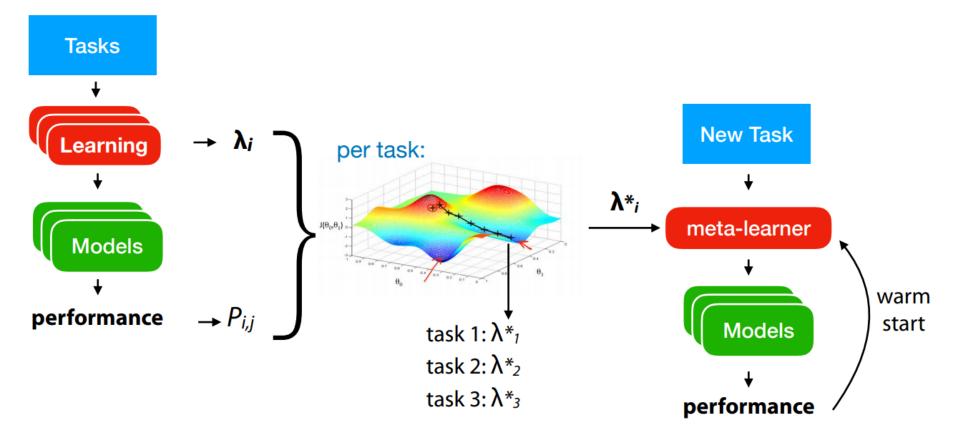
# **Top-K** recommendation

- Build a global (multi-objective) ranking, recommend the top-K
- Requires fixed selection of candidate configurations (portfolio)
- Can be used as a warm start for optimization techniques



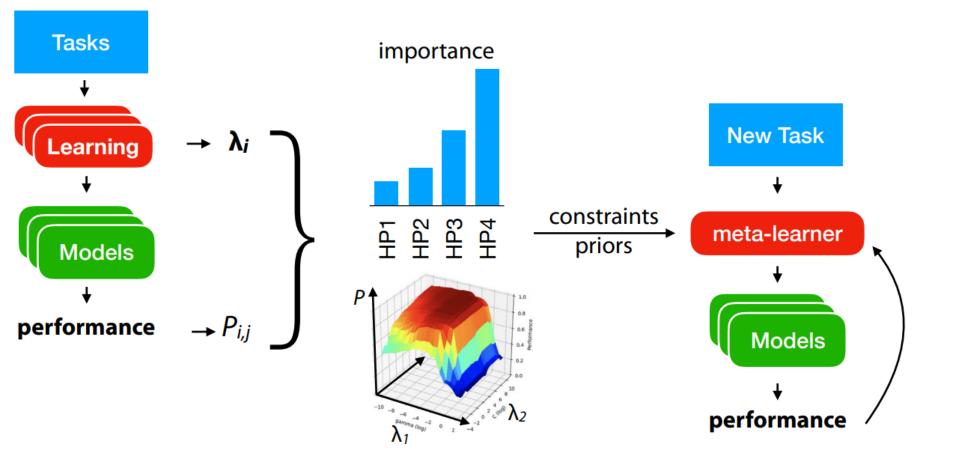
# Warm-starting with plugin estimators

- What if prior configurations are not optimal?
- Per task, fit a differentiable plugin estimator on all evaluated configurations
- Do gradient descent to find optimized configurations, recommend those



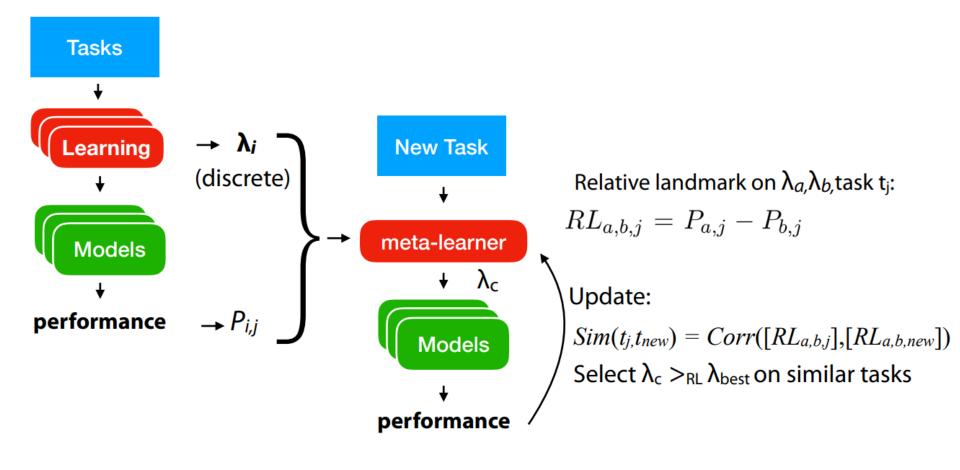
# Configuration space design

- Functional ANOVA: select hyperparameters that cause variance in the evaluations<sup>1</sup>
- **Tunability**: improvement from tuning a hyperparameter vs. using a good default<sup>2</sup>
- Search space pruning: exclude regions yielding bad performance on similar tasks<sup>3</sup>



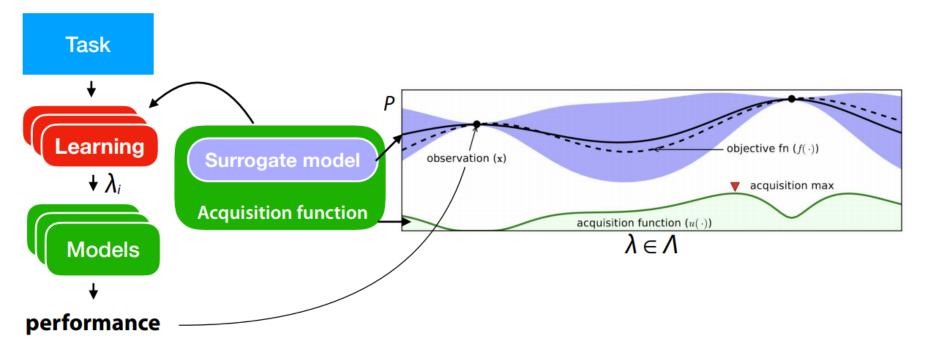
# Active testing

- Task are similar if observed relative performance of configurations is similar
- Tournament-style selection, warm-start with overall best configurations λ<sub>best</sub>
- Next candidate  $\lambda_c$ : the one that beats current  $\lambda_{best}$  on similar tasks (from portfolio)



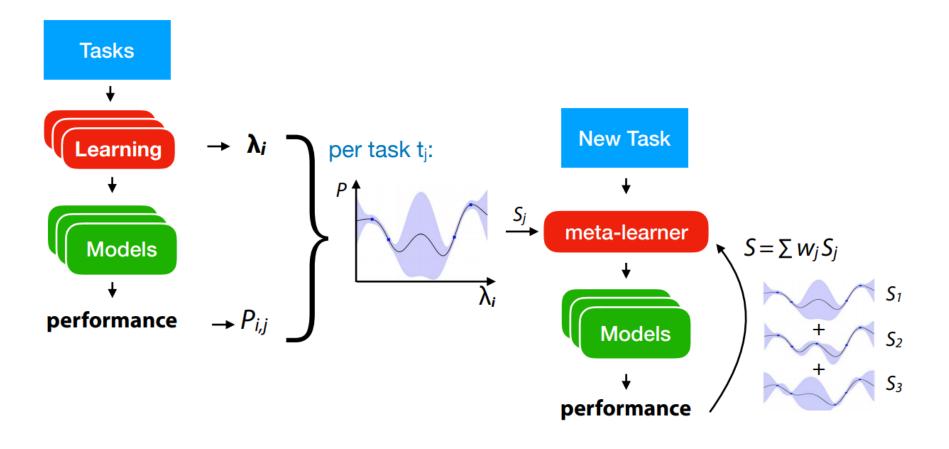
# Bayesian optimization (refresh)

- Learns how to learn within a single task (short-term memory)
- Surrogate model: *probabilistic* regression model of configuration performance
- Can we transfer what we learned to new tasks (long term memory)?



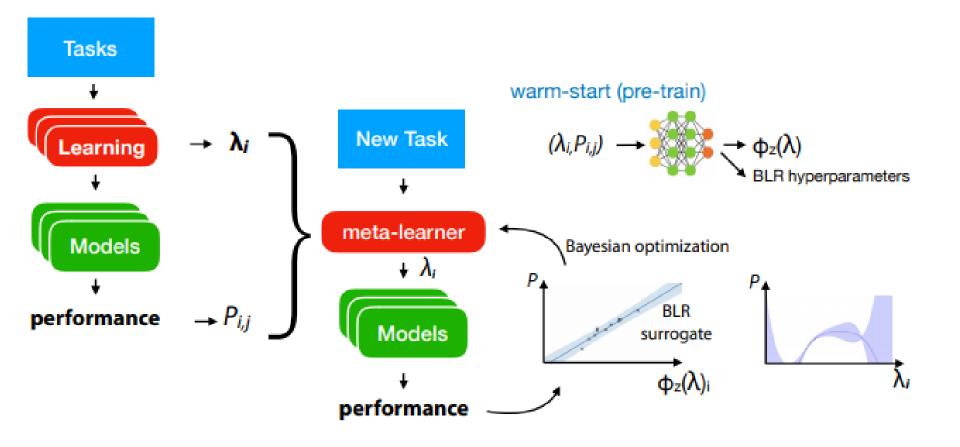
# Surrogate model transfer

- If task j is similar to the new task, its surrogate model S<sub>j</sub> will do well
- Sum up all S<sub>j</sub> predictions, weighted by task similarity (relative landmarks)<sup>1</sup>
- Build combined Gaussian process, weighted by current performance on new task<sup>2</sup>



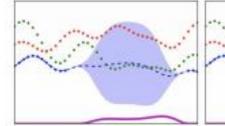
# Warm-started multi-task learning

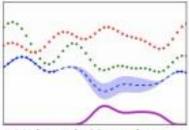
- Bayesian linear regression (BLR) surrogate model on every task
- Learn a suitable basis expansion  $\phi_z(\lambda)$ , joint representation for all tasks
- Scales linearly in # observations, transfers info on configuration space



# **Multi-task Bayesian optimization**

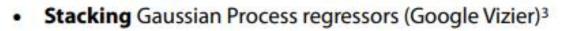
- Multi-task Gaussian processes: train surrogate model on t tasks simultaneously<sup>1</sup>
  - If tasks are similar: transfers useful info
  - Not very scalable



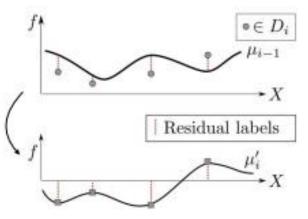


- Independent GP predictions
- Multi-task GP predictions

- Bayesian Neural Networks as surrogate model<sup>2</sup>
  - Multi-task, more scalable

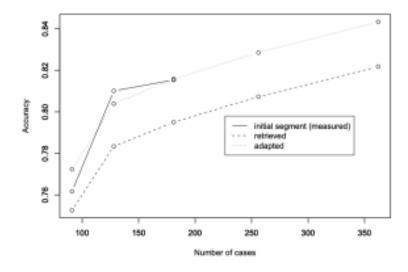


- Sequential tasks, each similar to the previous one
- Transfers a prior based on residuals of previous GP



## Other techniques

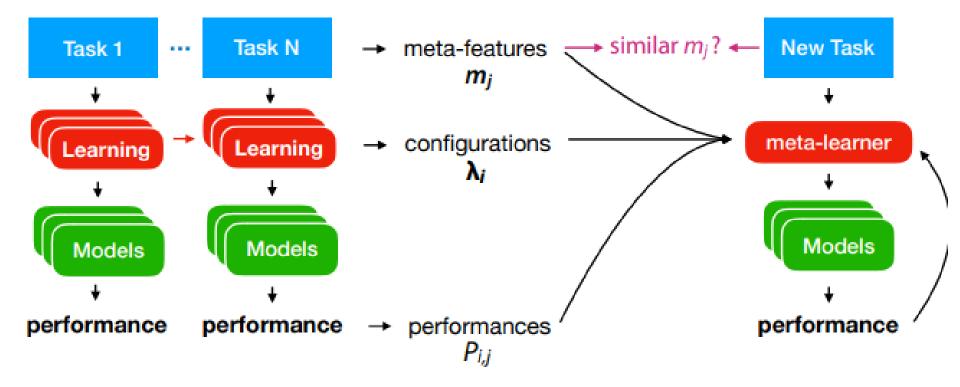
- Transfer learning with multi-armed bandits<sup>1</sup>
  - View every task as an arm, learn to `pull` observations from the most similar tasks
  - Reward: accuracy of configurations recommended based on these observations
- Transfer learning curves<sup>2,3</sup>
  - Learn a partial learning curve on a new task, find best matching earlier curves
  - Predict the most promising configurations based on earlier curves



### 2. Reason about model performance across tasks

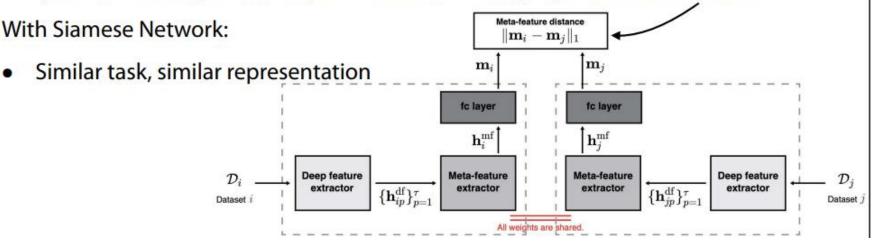
#### Meta-features: measurable properties of the tasks

(number of instances and features, class imbalance, feature skewness,...)



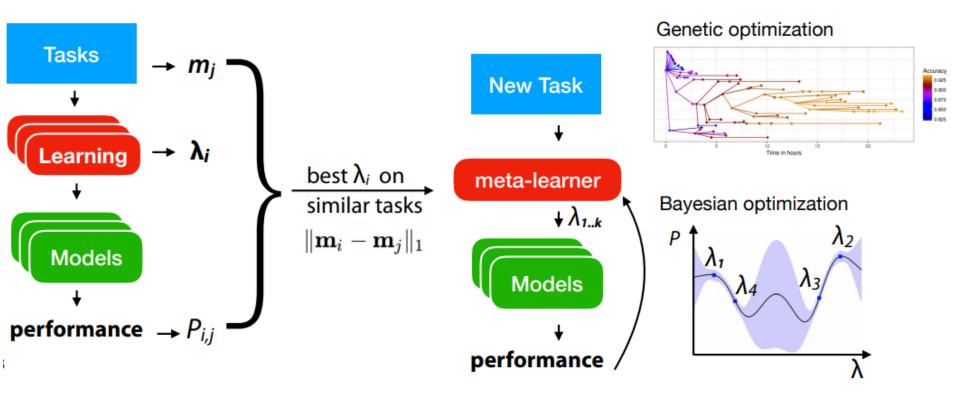
# Meta-features

- Hand-crafted (interpretable) meta-features<sup>1</sup>
  - Number of instances, features, classes, missing values, outliers,...
  - Statistical: skewness, kurtosis, correlation, covariance, sparsity, variance,...
  - Information-theoretic: class entropy, mutual information, noise-signal ratio,...
  - Model-based: properties of simple models trained on the task
  - Landmarkers: performance of fast algorithms trained on the task
  - Domain specific task properties
- Learning a joint task representation
  - Deep metric learning: learn a representation *h*<sup>mf</sup> using a ground truth distance<sup>2</sup>



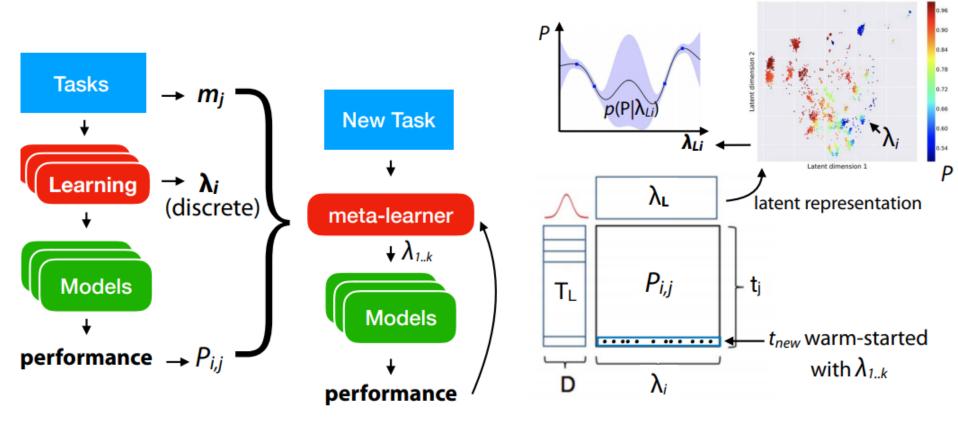
# Warm-starting from similar tasks

- Find k most similar tasks, warm-start search with best θ<sub>i</sub>
  - Genetic hyperparameter search<sup>1</sup>
  - Auto-sklearn: Bayesian optimization (SMAC)<sup>2</sup>
    - Scales well to high-dimensional configuration spaces



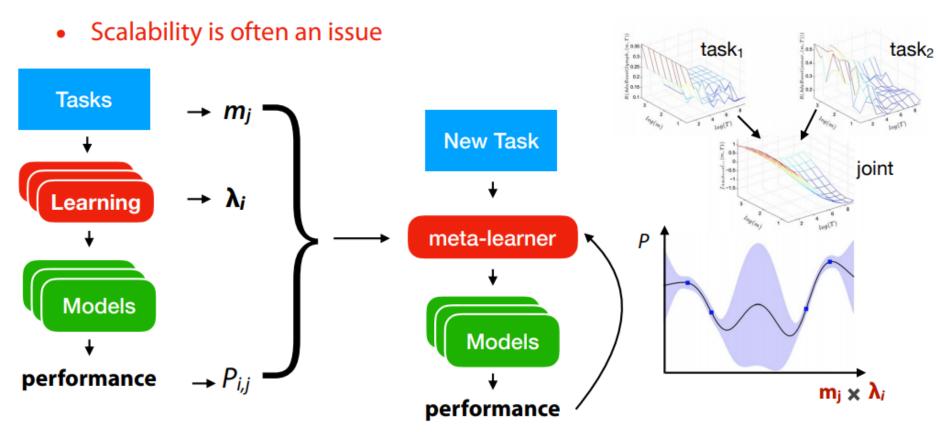
# Warm-starting from similar tasks

- Collaborative filtering: configurations λ<sub>i</sub> are `rated' by tasks t<sub>j</sub>
  - Probabilistic matrix factorization
    - Learns a latent representation for tasks and configurations
    - Returns probabilistic predictions for Bayesian optimization
    - Use meta-features to warm-start on new task



# Global surrogate models

- Train a task-independent surrogate model with meta-features in inputs
  - SCOT: Predict ranking of λ<sub>i</sub> with surrogate ranking model + m<sub>j</sub>.<sup>1</sup>
  - Predict P<sub>i,j</sub> with multilayer Perceptron surrogates + m<sub>j.</sub><sup>2</sup>
  - Build joint GP surrogate model on most similar ( $\|\mathbf{m}_i \mathbf{m}_j\|_2$ ) tasks. <sup>3</sup>



#### Meta-models

- Learn direct mapping between meta-features and Pi,j
  - Zero-shot meta-models: predict best λ<sub>i</sub> given meta-features <sup>1</sup>

 $m_j \rightarrow \text{meta-learner} \rightarrow \lambda_{best}$ 

• Ranking models: return ranking  $\lambda_{1..k}^2$ 

 $m_j \rightarrow \text{meta-learner} \rightarrow \lambda_{1..k}$ 

Predict which algorithms / configurations to consider / tune<sup>3</sup>

 $m_j \rightarrow \text{meta-learner} \rightarrow \Lambda$ 

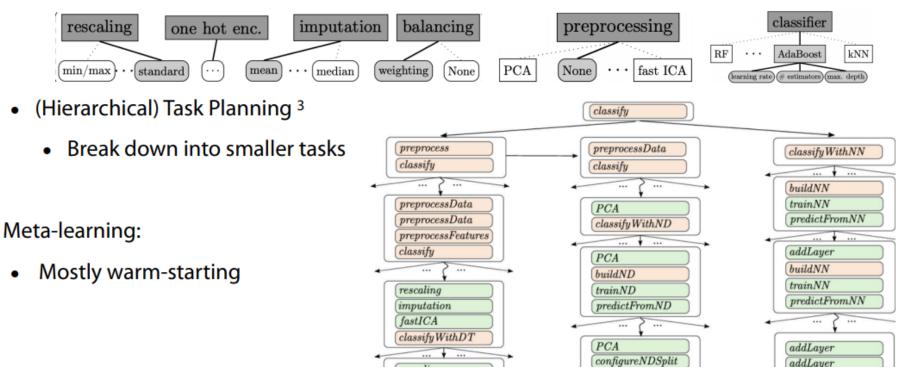
Predict performance / runtime for given θ<sub>i</sub> and task<sup>4</sup>

 $m_{j}, \lambda_{i} \rightarrow \text{meta-learner} \rightarrow P_{ij}$ 

Can be integrated in larger AutoML systems: warm start, guide search,...

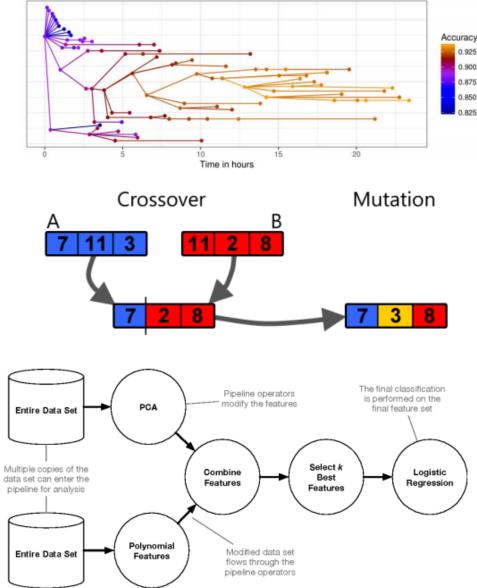
#### Learning Pipelines

- Compositionality: the learning process can be broken down into smaller tasks
  - Easier to learn, more transferable, more robust
- Pipelines are one way of doing this, but how to control the search space?
  - Select a fixed set of possible pipelines. Often works well (less overfitting)<sup>1</sup>
  - Impose a fixed structure on the pipeline <sup>2</sup>



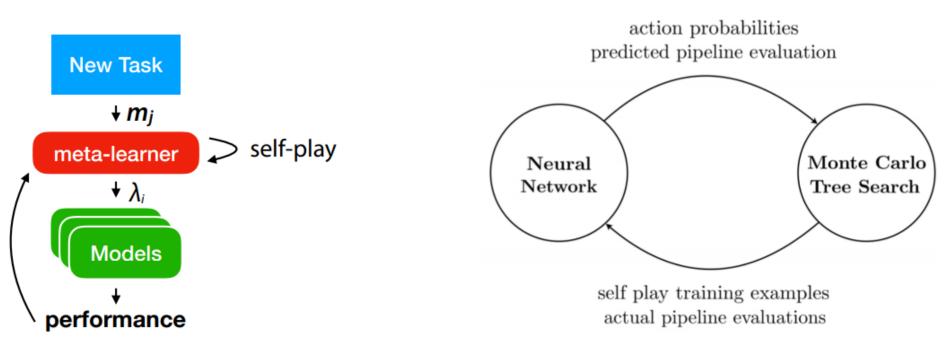
# **Evolving pipelines**

- Start from simple pipelines
- Evolve more complex ones if needed
- Reuse pipelines that do specific things
- Mechanisms:
  - Cross-over: reuse partial pipelines
  - Mutation: change structure, tuning
- Approaches:
  - TPOT: Tree-based pipelines<sup>1</sup>
  - GAMA: asynchronous evolution<sup>2</sup>
  - RECIPE: grammar-based<sup>3</sup>
- Meta-learning:
  - Largely unexplored
  - Warm-starting, meta-models



## Learning to learn through self-play

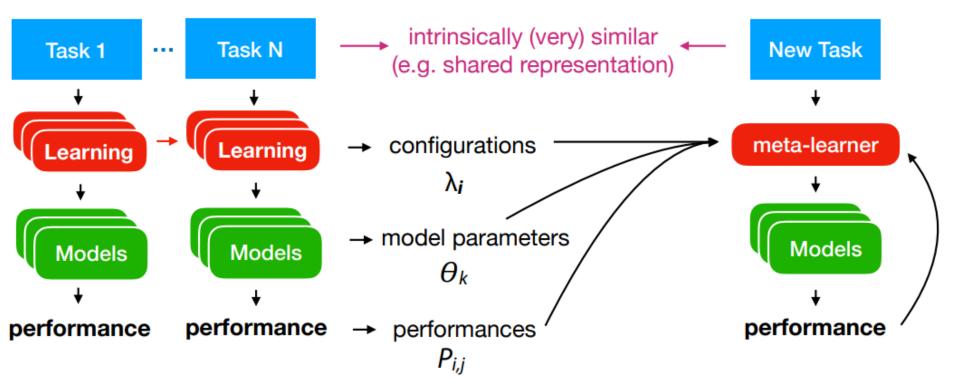
- Build pipelines by selecting among actions
  - insert, delete, replace pipeline parts
- Neural network (LSTM) receives task meta-features, pipelines and evaluations
  - Predict pipeline performance and action probabilities
- Monte Carlo Tree Search builds pipelines based on probabilities
  - Runs multiple simulations to search for a better pipeline



## 3. Learning from trained models

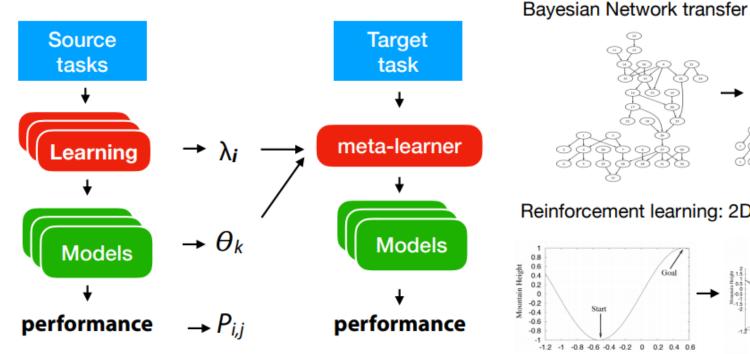
#### Models trained on similar tasks

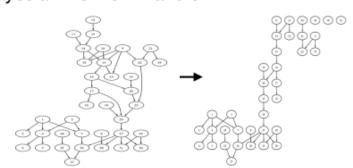
(model parameters, features,...)



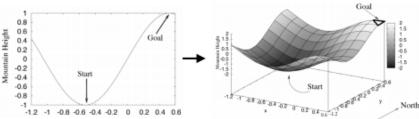
# Transfer Learning

- Select source tasks, transfer trained models to similar target task <sup>1</sup>
- Use as starting point for tuning, or *freeze* certain aspects (e.g. structure)
  - Bayesian networks: start structure search from prior model <sup>2</sup>
  - Reinforcement learning: start policy search from prior policy <sup>3</sup>



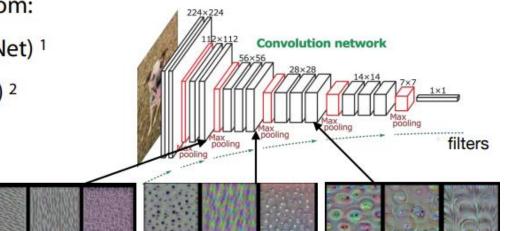


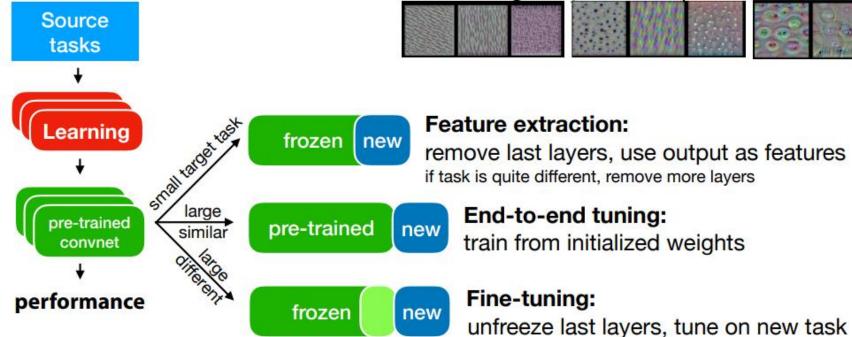
Reinforcement learning: 2D to 3D mountain car



## Transfer features, initializations

- For neural networks, both structure and weights can be transferred
- Features and initializations learned from:
  - Large image datasets (e.g. ImageNet)<sup>1</sup>
  - Large text corpora (e.g. Wikipedia)<sup>2</sup>
- Fails if tasks are not similar enough <sup>3</sup>

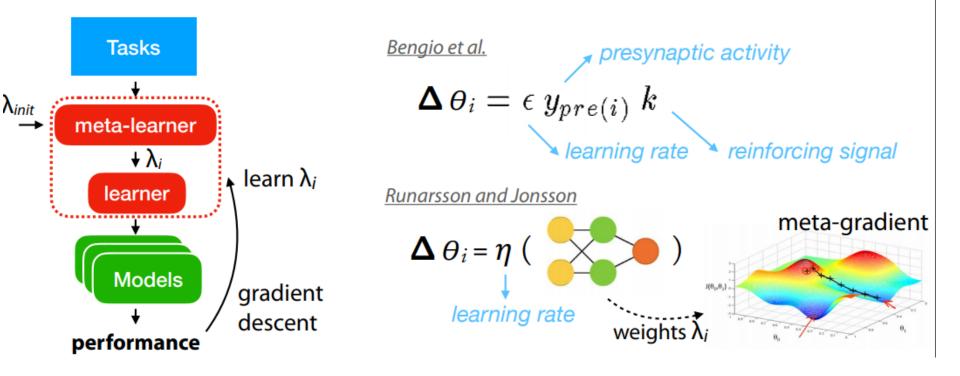




#### Learning to learn by gradient descent

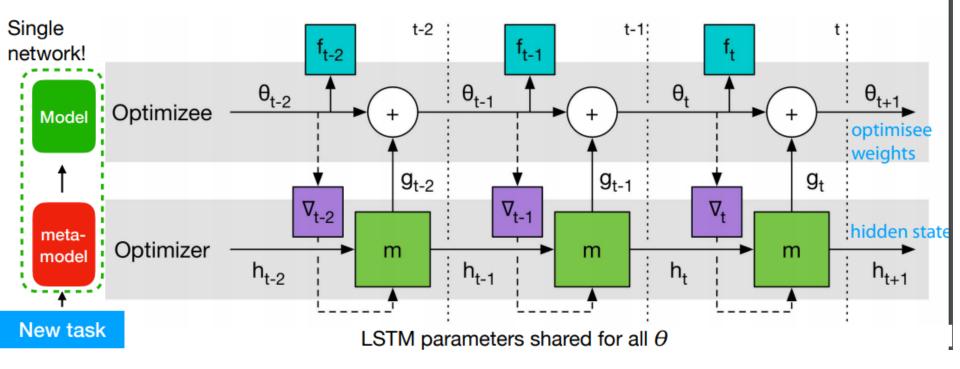
- Our brains *probably* don't do backprop, replace it with:
  - Simple parametric (bio-inspired) rule to update weights <sup>1</sup>
  - Single-layer neural network to learn weight updates <sup>2</sup>
- Learn parameters across tasks, by gradient descent (meta-gradient)

 $\Delta w_{ii}$ 



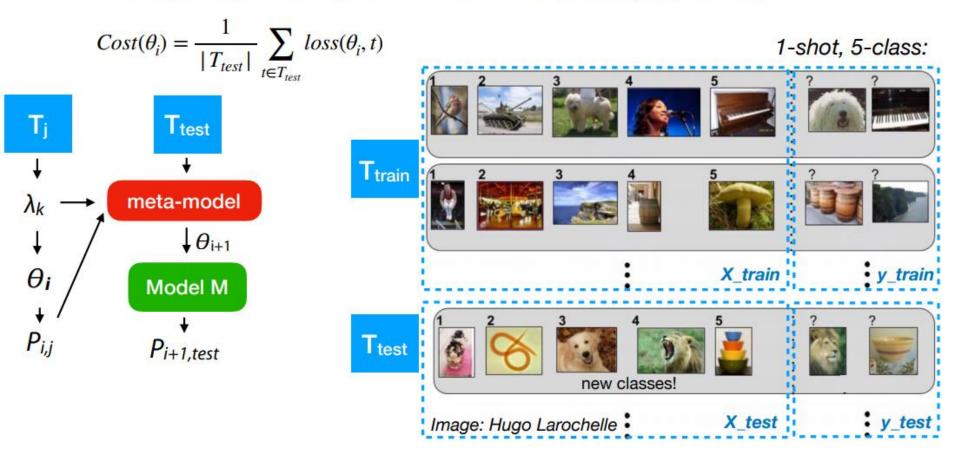
#### Learning to learn gradient descent by gradient descent

- Replace backprop with a recurrent neural net (LSTM)<sup>1</sup>, not so scalable
- Use a coordinatewise LSTM [m] for scalability/flexibility (cfr. ADAM, RMSprop)<sup>2</sup>
  - Optimizee: receives weight update g<sub>t</sub> from optimizer
  - Optimizer: receives gradient estimate  $\nabla_t$  from optimizee
    - Learns how to do gradient descent across tasks



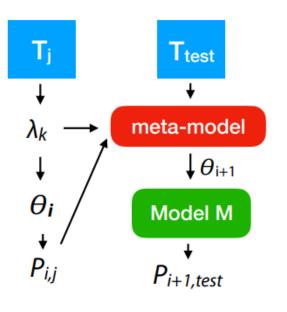
## Few-shot learning

- Learn how to learn from few examples (given similar tasks)
  - Meta-learner must learn how to train a base-learner based on prior experience
    - Parameterize base-learner model and learn the parameters  $\theta_i$



## Few-shot learning: approaches

. . .



- LSTM + gradient descent
   <u>Ravi and Larochelle 2017</u>
- Learn  $\theta_{init+}$  gradient descent
- kNN-like: Memory + similarity
- Learn embedding + classifier

<u>Finn et al. 2017</u>

Vinyals et al. 2016

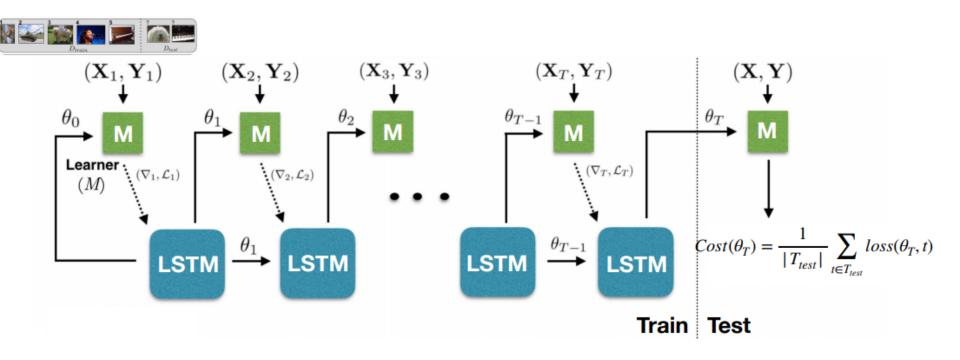
<u>Snell et al. 2017</u>

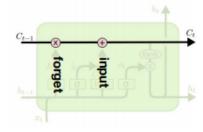
- $Cost(\theta_i) = \frac{1}{|T_{test}|} \sum_{t \in T_{test}} loss(\theta_i, t)$
- Black-box meta-learner
  - Neural Turing machine (with memory) Santoro et al. 2016
  - Neural attentive learner

- Mishus at al 2010
- <u>Mishra et al. 2018</u>

#### LSTM meta-learner + gradient descent

- Gradient descent update  $\theta_t$  is similar to LSTM cell state update  $c_t$ 
  - $\theta_t = \theta_{t-1} \alpha_t \nabla_{\theta_{t-1}} \mathcal{L}_t \qquad c_t = f_t \odot c_{t-1} + i_t \odot \tilde{c}_t$
- Hence, training a meta-learner LSTM yields an update rule for training M
  - Start from initial  $\theta_{0}$ , train model on first batch, get gradient and loss update
  - Predict  $\theta_{t+1}$ , continue to t=T, get cost, backpropagate to learn LSTM weights, optimal  $\theta_0$





# Model-agnostic meta-learning

- Quickly learn new skills by learning a model *initialization* that generalizes better to similar tasks
   — meta-learning
  - Current initialization  $\theta$
  - On K examples/task, evaluate  $\nabla_{\theta} L_{T_i}(f_{\theta})$
  - Update weights for  $\theta_1, \theta_2, \theta_3$
  - Update θ to minimize sum of per-task losses
  - Repeat  $\theta \leftarrow \theta \beta \nabla_{\theta} \sum_{i=1}^{n}$

task losses  

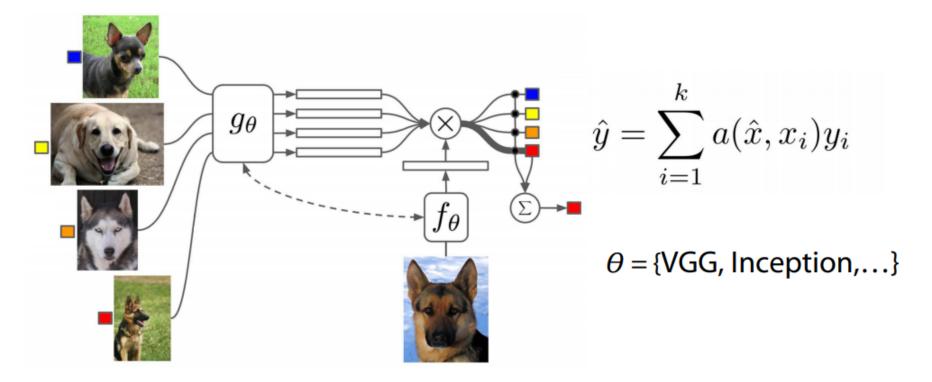
$$\nabla \mathcal{L}_{1}$$
 $\nabla \mathcal{L}_{2}$ 
 $\nabla \mathcal{L}_{1}$ 
 $\theta_{1}^{*}$ 

---- learning/adaptation

- More resilient to overfitting
- Generalizes better than LSTM approaches
- Universality: no theoretical downsides in terms of expressivity when compared to alternative meta-learning models.
- REPTILE: do SGD for k steps in one task, only then update initialization weights<sup>3</sup>

#### 1-shot learning with Matching networks

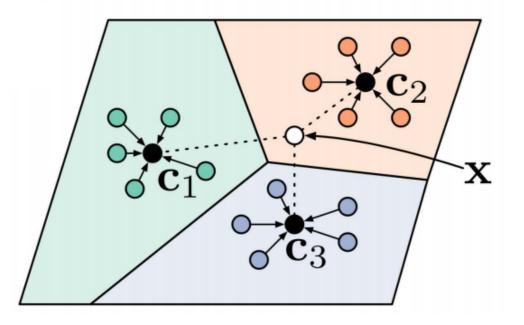
- Don't learn model parameters, use non-parameters model (like kNN)
- Choose an embedding network f and g (possibly equal)
- Choose an attention kernel  $a(\hat{x}, x_i)$ , e.g. softmax over cosine distance
- Train complete network in minibatches with few examples per task



# Prototypical networks

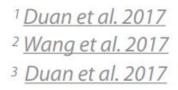
<u>Snell et al. 2017</u> <u>Ren et al. 2018</u>

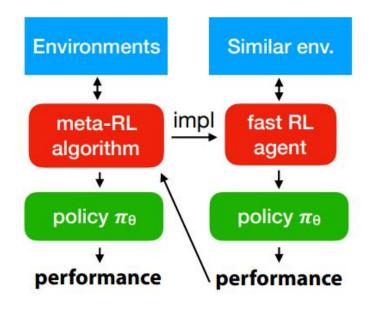
- Train a "prototype extractor" network
- Map examples to p-dimensional embedding so examples of a given class are close together
- Calculate a prototype (mean vector) for every class
- Map test instances to the same embedding, use softmax over distance to prototype
- Using more classes during meta-training works better!

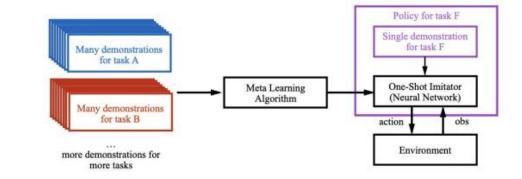


## Learning to reinforcement learn

- Humans often learn to play new games much faster than RL techniques do
- Reinforcement learning is very suited for learning-to-learn:
  - Build a learner, then use performance as that learner as a reward
- Learning to reinforcement learn 1,2
  - Use RNN-based deep RL to train a recurrent network on many tasks
  - · Learns to implement a 'fast' RL agent, encoded in its weights







- Also works for few-shot learning <sup>3</sup>
  - Condition on observation + upcoming demonstration
- You don't know what someone is trying to teach you, but you prepare for the lesson

#### Learning to learn more tasks

- Active learning
  - Deep network (learns representation) + policy network
  - Receives state and reward, says which points to query next
- Density estimation
  - Learn distribution over small set of images, can generate new ones
  - Uses a MAML-based few-shot learner
- Matrix factorization
  - Deep learning architecture that makes recommendations
  - Meta-learner learns how to adjust biases for each user (task)
- Replace hand-crafted algorithms by learned ones.
- Look at problems through a meta-learning lens!

Pang et al. 2018

Reed et al. 2017

<u>Vartak et al. 2017</u>

## Meta-data sharing building a shared memory Vanschoren et al. 2014

- OK, but how do I get large amounts of meta-data for meta-learning?
- OpenML.org
  - Thousands of uniform datasets
  - 100+ meta-features
  - Millions of evaluated runs
    - Same splits, 30+ metrics
    - Traces, models (opt)
  - APIs in Python, R, Java,...
  - Publish your own runs
  - Never ending learning
  - Benchmarks

Open positions! Scientific programmer Teaching PhD

```
import openml as oml
from sklearn import tree
task = oml.tasks.get_task(14951)
clf = tree.ExtraTreeClassifier()
flow = oml.flows.sklearn_to_flow(clf)
run = oml.runs.run_flow_on_task(task, flow)
myrun = run.publish()
```

run locally, share globally



Joaquin Vanschoren **Hilde Weerts** edorigatti frontier . loel Goossens Niels Hellinga Minopeivu Zhano Evertian Peer stevens jethefer Hongliang Qiu . Yezi Zhu János Szedelényi Chin-Fang Lin Wenting Xion Lirong Zhang M de Roode Tianyu Zhou **Ruud Andriessen** Angelo Majoor

# Towards human-like learning to learn

- Learning-to-learn gives humans a significant advantage
  - Learning how to learn any task empowers us far beyond knowing how to learn specific tasks.
  - It is a **universal** aspect of life, and how it evolves
- Very exciting field with many unexplored possibilities
  - Many aspects not understood (e.g. task similarity), need more experiments.
- Challenge:
  - Build learners that never stop learning, that learn from each other
  - Build a *global memory* for learning systems to learn from
  - Let them explore by themselves, active learning